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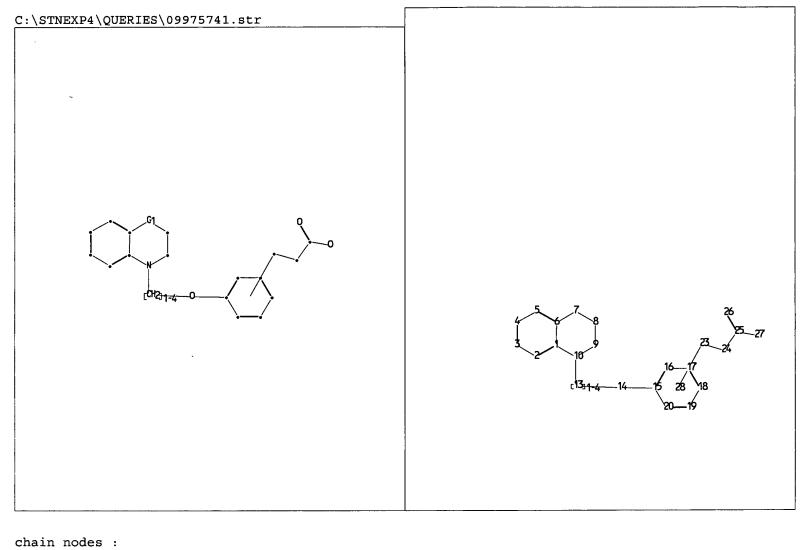
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```
13 14 23 24 25 26 27
ring nodes :
    1 2 3 4 5 6 7 8 9 10 15 16 17 18 19 20
chain bonds :
    10-13 13-14 14-15 23-24 24-25 25-26 25-27
ring bonds :
    1-2 1-6 1-10 2-3 3-4 4-5 5-6 6-7 7-8 8-9 9-10 15-16 15-20 16-17 17-18 18-19 19-20
exact/norm bonds :
    1-10 6-7 7-8 8-9 9-10 10-13 13-14 14-15 23-24 24-25 25-26 25-27
normalized bonds :
    1-2 1-6 2-3 3-4 4-5 5-6 15-16 15-20 16-17 17-18 18-19 19-20
isolated ring systems :
    containing 1 : 15 :
```

G1:0,S

Match level:
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 13:CLASS
14:CLASS 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 23:CLASS 24:CLASS
25:CLASS 26:CLASS 27:CLASS 28:CLASS

=>

Uploading 09975741.str

L1 STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1 STF

G1 0, S

Structure attributes must be viewed using STN Express query preparation.

=> s 11 sss sam

SAMPLE SEARCH INITIATED 15:25:11 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 59 TO ITERATE

100.0% PROCESSED 59 ITERATIONS 11 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 720 TO 1640

PROJECTED ANSWERS: 22 TO 418

L2 11 SEA SSS SAM L1

=> s l1 sss ful

FULL SEARCH INITIATED 15:26:09 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1089 TO ITERATE

100.0% PROCESSED 1089 ITERATIONS 171 ANSWERS

SEARCH TIME: 00.00.01

L3 171 SEA SSS FUL L1

=> s 13

L4 5 L3

=> d 14 1-5 bib, ab, hitstr

```
L4
     ANSWER 1 OF 5 CAPLUS COPYRIGHT 2003 ACS on STN
      2003:633683 CAPLUS
AN
DN
     139:185673
ΤI
      Preparation and compositions of polymorphic forms of bicyclic antidiabetic
      agents
IN
      Srisilla, Raju; Potlapally, Rajender Kumar; Mamillapalli, Ramabhadra
      Sarma; Gaddam, Om Reddy
     Reddy's Laboratories Limited, India
PA
SO
     PCT Int. Appl., 35 pp.
     CODEN: PIXXD2
DT
      Patent
LΑ
     English
FAN.CNT 1
     PATENT NO.
                         KIND
                                ΊDATE
                                                 APPLICATION NO.
                                                                     DATE
                                20030814
                          A1
                                                 WO 2003-IB408
                                                                     20030207
PΙ
     WO 2003066612
                                          AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
              AE, AG, AL, AM, AT, AU,
              CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD,
              RU, TJ, TM
          RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG,
              CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC,
              NL, PT, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW,
              ML, MR, NE, SN, TD, TG
PRAI IN 2002-MA95
                                20020207
                          Α
     This invention relates to novel polymorphic/pseudopolymorphic forms and
     compns. of arginine salt of 3-[4-[2-(3,4-dihydro-1,4-benzothiazin-4-
     yl)ethoxy]phenyl]-2-ethoxypropanoic acid, preferably, L-arginine salt of
      (2S)-3-[4-[2-(3,4-dihydro-1,4-benzothiazin-4-yl)ethoxy]phenyl]-2-
     ethoxypropanoic acid. The polymorphic forms of the present invention are
     more active, as antidiabetic and hypolipidemic agent, than the
     3-[4-[2-(2,3-dihydro-1,4-benzothiazin-4-yl)ethoxy]phenyl]-2-
     ethoxypropanoic acid.
ΙT
     305827-56-9P
     RL: PEP (Physical, engineering or chemical process); PNU (Preparation,
     unclassified); PRP (Properties); PYP (Physical process); THU (Therapeutic
     use); BIOL (Biological study); PREP (Preparation); PROC (Process); USES
         (prepn. and compns. of polymorphic forms of arginine salt of
         benzothiazine antidiabetic/hypolipidemic agent)
RN
     305827-56-9 CAPLUS
     L-Arginine, mono[(.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-
CN
     yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)
     CM
           1
     CRN
           223607-27-0
     CMF
           C21 H25 N O4 S
```

CRN 74-79-3 CMF C6 H14 N4 O2

Absolute stereochemistry.

IT 223607-27-0

RL: RCT (Reactant); RACT (Reactant or reagent)
 (prepn. and compns. of polymorphic forms of arginine salt of
 benzothiazine antidiabetic/hypolipidemic agent)

RN 223607-27-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)- (9CI) (CA INDEX NAME)

```
ANSWER 2 OF 5 CAPLUS COPYRIGHT 2003 ACS on STN
L4
ΑN
     2003:242313 CAPLUS
DN
     138:255240
TI
     An improved process for the preparation of bicyclic antidiabetic
     compounds, specifically (S)-3-[4-[2-(3,4-dihydro-2H-benzo[b][1,4]thiazin-4-
     yl)ethoxy]phenyl]-2-alkoxypropanoic acids and their oxazin-4-yl analogs,
     using optimized bases and solvents for etherification
     Batchu, Chandrasekhar; Addanki, Sivaramaprasad; Mamillapalli,
TN
     Ramabhadrasarma; Gaddam, Omreddy
     Reddy's Laboratories Ltd., India
PA
SO
     PCT Int. Appl., 14 pp.
     CODEN: PIXXD2
DT
     Patent
LΑ
     English
FAN.CNT 1
     PATENT NO.
                       KIND
                              DATE
                                             APPLICATION NO.
                                                               DATE
PΙ
     WO 2003024944
                        A1
                             20030327
                                             WO 2002-IB3875
                                                               20020920
             AE, AG, AL, AM, AĮ, AU,
                                      AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
             CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,
             GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
             LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD,
             RU, TJ, TM
         RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG,
             CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL,
             PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR,
             NE, SN, TD, TG
PRAI IN 2001-MA778
                             20010920
                        Α
     CASREACT 138:255240; MARPAT 138:255240
os
AΒ
     The invention relates to an improved process for the prepn. of
     antidiabetic compds. I [wherein: R1 = (C1-C6)alkyl; X = S or O atom].
     method involves (1) etherification of phenol derivs. II [R1, R2 =
     (C1-C6)alkyl] with mesylates III in the presence of a base and solvent at
     10-150.degree. over 2-30 h to give esters IV; (2) hydrolysis of IV by a
     base and a solvent at 10-40.degree. over 20 min to 12 h; and (3) isolation
     of I by conventional methods. In particular, the base in both steps is
     selected from Na2CO3, K2CO3, NaOH, or KOH, and the solvent is selected
     from toluene, xylene, MeOH, EtOH, n-PrOH, iso-PrOH, or their mixts. I are
     known antidiabetic agents, with prepns. previously described in US
     6,265,401 and WO 00/66572. Disadvantages of these prior art methods
     include moisture-sensitive reactions, use of column chromatog., low
     yields, long cycle times, impure products, and an inconvenient solvent
     (DMF). The new method is simple, safe, scalable to com. prodn., and uses
     inexpensive chems. The new solvents give improved yields and drastically
     reduced cycle times, making the process com. viable. For instance,
     etherification of (-)-II [R1 = R2 = Et] with III [X = S] using K2CO3 in
     refluxing toluene for 4-5 h with azeotropic removal of H2O gave 92% (-)-IV
     [R1 = R2 = Et; X = S] of 97% purity. The latter ester in MeOH was
     hydrolyzed by 10% NaOH soln. at <20.degree. and then at 25-35.degree.,
     followed by extn. with PhMe and acidification with 25% H2SO4, to give 82%
     (-)-I [R1 = R2 = Et; X = S], also of 97% purity. Nearly identical results
     were obtained for the case of R1 = Et, R2 = iso-Pr, X = S.
IT
     305827-41-2P, (S)-(-)-Ethyl 3-[4-[2-(3,4-dihydro-2H-
     benzo[b][1,4]thiazin-4-yl)ethoxy]phenyl]-2-ethoxypropanoate
     503056-17-5P, (S)-(-)-Isopropyl 3-[4-[2-(3,4-dihydro-2H-
     benzo[b][1,4]thiazin-4-yl)ethoxy]phenyl]-2-ethoxypropanoate
```

RL: IMF (Industrial manufacture); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent) (intermediate; improved prepn. of bicyclic antidiabetic compds. by etherification of phenol and mesylate derivs.)

RN 305827-41-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy].alpha.-ethoxy-, ethyl ester, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 503056-17-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, 1-methylethyl ester, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 223607-27-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RE.CNT 1 THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

```
L4
     ANSWER 3 OF 5 CAPLUS COPYRIGHT 2003 ACS on STN
     2002:293636 CAPLUS
AN
DN
     136:319405
ΤI
     Salts of benzothiazine and benzoxazine derivatives and pharmaceutical
     compositions containing them
TN
     Gaddam, Om Reddy; Batchu, Chandrasekhar; Mamillapalli, Ramabhadra Sarma;
     Addanki, Siva Rama Prasad
     Reddy's Research Foundation, India
PA
SO
     PCT Int. Appl., 58 pp.
     CODEN: PIXXD2
DT
     Patent
LA
     English
FAN.CNT 1
     PATENT NO.
                       KIND DATE
                                              APPLICATION NO.
                                                                 DATE
     _____
                       ____
     WO 2002030914
                              20020418
                                              WO 2001-IN175
                                                                 20011010
PΙ
                        A1
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              CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,
              GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
              LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PH, PL,
         PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF,
              BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
     AU 2002012672
                        A5
                              20020422
                                              AU 2002-12672
                                                                 20011010
     US 2002142993
                        A1
                              20021003
                                              US 2001-975741
                                                                 20011010
                              20001012
PRAI IN 2000-MA863
                        Α
     US 2001-266996P
                        Ρ
                              20010205
     WO 2001-IN175
                        W
                              20011010
os
     MARPAT 136:319405
     The present invention relates to pharmaceutically acceptable salts of
AB
     compd. of the general formula (I), their derivs., their analogs, their
     tautomeric forms, their stereoisomers, their polymorphs, their
     pharmaceutically acceptable solvates and pharmaceutically acceptable
     compns. contg. them. wherein R1 represents hydrogen, halogen atom such as
     fluorine, chlorine, bromine or iodine; hydroxy, nitro, cyano or lower
     alkyl group; R2 represents hydrogen, lower alkyl or oxo group; X
     represents a heteroatom selected from oxygen or sulfur; R3 represents
     hydrogen or lower alkyl group; n is an integer ranging from 1-4; M
     represents a counter ion or a moiety which forms a pharmaceutically
     acceptable salt; p is an integer ranging from 1 to 2. Examples are
     provided showing effectiveness of such compds. in inhibiting HMG CoA
     reductase activity, and in lowering plasma triglyceride and cholesterol
     levels.
IT
     412316-40-6P
     RL: PAC (Pharmacological activity); PKT (Pharmacokinetics); SPN (Synthetic
     preparation); THU (Therapeutic use); BIOL (Biological study); PREP
     (Preparation); USES (Uses)
         (salts of benzothiazines and benzoxazines for treatment of various
        diseases such as atherosclerosis and diabetes)
RN
     412316-40-6 CAPLUS
CN
     Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-
```

diamide (1:1) (9CI) (CA INDEX NAME)

.alpha.-ethoxy-, (.alpha.S)-, compd. with N,N-dimethylimidodicarbonimidic

CRN 223607-27-0 CMF C21 H25 N O4 S

Absolute stereochemistry. Rotation (-).

chitia peries

CM 2

CRN 657-24-9 CMF C4 H11 N5

NH NH || || Me<sub>2</sub>N-C-NH-C-NH<sub>2</sub>

IT 412316-39-3P 412316-41-7P 412316-42-8P 412316-43-9P 412316-44-0P 412316-45-1P 412316-47-3P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(salts of benzothiazines and benzoxazines for treatment of various diseases such as atherosclerosis and diabetes)

RN 412316-39-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with (.beta.S)-.beta.-aminobenzeneethanol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0 CMF C21 H25 N O4 S

CRN 20989-17-7 CMF C8 H11 N O

Absolute stereochemistry. Rotation (+).

$$\begin{array}{c|c} & & \text{NH2} \\ & & \\ &$$

RN · 412316-41-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with N-cyclohexylcyclohexanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0 CMF C21 H25 N O4 S

CRN 101-83-7 CMF C12 H23 N

RN 412316-42-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with (.alpha.R)-.alpha.-methylbenzenemethanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0 CMF C21 H25 N O4 S

Absolute stereochemistry. Rotation (-).

CM 2

CRN 3886-69-9 CMF C8 H11 N

Absolute stereochemistry.

RN 412316-43-9 CAPLUS

CN L-Lysine, mono[(.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CRN 223607-27-0 CMF C21 H25 N O4 S

Absolute stereochemistry. Rotation (-).

CM 2

CRN 56-87-1

CMF C6 H14 N2 O2

Absolute stereochemistry.

RN 412316-44-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with hydrazinecarboximidamide (1:1), monoformate (9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0

CMF C21 H25 N O4 S

CRN 79-17-4 C H6 N4 CMF

NH H<sub>2</sub>N- "С- NH- NH<sub>2</sub>

> CM3

CRN 64-18-6 CMF C H2 O2

o = ch - oh

RN 412316-45-1 CAPLUS

Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-CN .alpha.-ethoxy-, (.alpha.S)-, compd. with 2-amino-2-(hydroxymethyl)-1,3propanediol (1:1) (9CI) (CA INDEX NAME)

CM

CRN 223607-27-0

CMF C21 H25 N O4 S

CRN 77-86-1 CMF C4 H11 N O3

$$\begin{array}{c} & \text{NH}_2 \\ | \\ \text{HO-CH}_2 - \text{C-CH}_2 - \text{OH} \\ | \\ \text{CH}_2 - \text{OH} \end{array}$$

RN 412316-47-3 CAPLUS

CN D-Glucitol, 1-deoxy-1-(methylamino)-, (.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0 . CMF C21 H25 N O4 S

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CM
      2
```

6284-40-8 CRN CMF C7 H17 N O5

Absolute stereochemistry.

```
OH
MeNH
         ŌН
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IT
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     412316-51-9 412316-52-0 412316-53-1
     412316-54-2 412316-55-3 412316-56-4
     412316-57-5 412316-58-6 412316-59-7
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     412316-66-6 412316-67-7 412316-68-8
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     412317-73-8 412317-74-9 412317-76-1
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     412915-59-4 412915-60-7 412915-63-0
     412915-66-3 412915-68-5 412915-69-6
     412915-70-9
    RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
        (salts of benzothiazines and benzoxazines for treatment of various
        diseases such as atherosclerosis and diabetes)
RN
    412316-48-4 CAPLUS
CN
    L-Lysine, mono[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-
    ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)
```

CM

223607-15-6 CRN CMF C21 H25 N O5

CRN 56-87-1

CMF C6 H14 N2 O2

Absolute stereochemistry.

RN 412316-49-5 CAPLUS

CN L-Lysine, mono[(.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8

CMF C21 H25 N O5

CRN 56-87-1

CMF C6 H14 N2 O2

Absolute stereochemistry.

RN 412316-50-8 CAPLUS

CN L-Lysine, mono[4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7

CMF C21 H25 N O4 S

CRN 56-87-1 CMF C6 H14 N2 O2

Absolute stereochemistry.

RN 412316-51-9 CAPLUS

CN L-Lysine, mono[(.alpha.R)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 56-87-1

CMF C6 H14 N2 O2

Absolute stereochemistry.

RN 412316-52-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with N-cyclohexylcyclohexanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CRN 101-83-7 CMF C12 H23 N

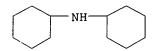
RN 412316-53-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with N-cyclohexylcyclohexanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 101-83-7 CMF C12 H23 N



RN 412316-54-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with N-cyclohexylcyclohexanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

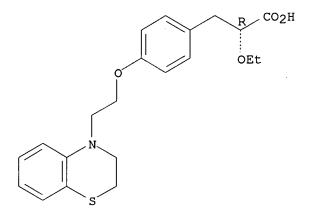
CRN 101-83-7 CMF C12 H23 N

RN 412316-55-3 CAPLUS

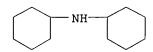
CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, compd. with N-cyclohexylcyclohexanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S



CRN 101-83-7 CMF C12 H23 N



RN 412316-56-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-ethoxy-, compd. with N,N-dimethylimidodicarbonimidic diamide (1:1)
(9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CRN 657-24-9 CMF C4 H11 N5

RN 412316-57-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with N,N-dimethylimidodicarbonimidic diamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 657-24-9 CMF C4 H11 N5

RN 412316-58-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy].alpha.-ethoxy-, compd. with N,N-dimethylimidodicarbonimidic diamide (1:1)
(9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 657-24-9 CMF C4 H11 N5

$$\begin{array}{ccccc} & \text{NH} & \text{NH} \\ & || & & || \\ \text{Me}_2 \text{N-C-NH-C-NH}_2 \end{array}$$

RN 412316-59-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, compd. with N,N-dimethylimidodicarbonimidic diamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 657-24-9 CMF C4 H11 N5

$$\begin{array}{c|c} & \text{NH} & \text{NH} \\ & || & || \\ \text{Me}_2 \text{N-C-NH-C-NH}_2 \end{array}$$

RN 412316-60-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with hydrazinecarboximidamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 79-17-4 CMF C H6 N4

ин || н<sub>2</sub>и-с-ин-ин<sub>2</sub>

RN 412316-61-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy].alpha.-ethoxy-, (.alpha.S)-, compd. with hydrazinecarboximidamide (1:1)
(9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0 CMF C21 H25 N O4 S

CRN 79-17-4 CMF C H6 N4

RN 412316-62-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, compd. with hydrazinecarboximidamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 79-17-4 CMF C H6 N4

 $\begin{array}{c} \text{nh} \\ || \\ \text{H}_2 \text{n-c-nh-nh}_2 \end{array}$ 

RN 412316-63-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with hydrazinecarboximidamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CM 2

CRN 79-17-4 CMF C H6 N4

ин || н2и-с-ин-ин2

RN 412316-64-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with hydrazinecarboximidamide (1:1)

(9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

Absolute stereochemistry. Rotation (-).

CM 2

CRN 79-17-4 CMF C H6 N4

ин || н<sub>2</sub>и- с- ин- ин<sub>2</sub>

RN 412316-65-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with (.alpha.R)-.alpha.-methylbenzenemethanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CRN 3886-69-9 CMF C8 H11 N

Absolute stereochemistry.

RN 412316-66-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with (.alpha.R)-.alpha.-methylbenzenemethanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 3886-69-9 CMF C8 H11 N

Absolute stereochemistry.

RN 412316-67-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with (.alpha.R)-.alpha.-methylbenzenemethanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 3886-69-9 CMF C8 H11 N

Absolute stereochemistry.

RN 412316-68-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, compd. with (.alpha.R)-.alpha.-methylbenzenemethanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 3886-69-9 CMF C8 H11 N

Absolute stereochemistry.

RN 412316-69-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with hydrazinecarboximidamide (1:1), monoformate (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 79-17-4 CMF C H6 N4

CM 3

CRN 64-18-6 CMF C H2 O2

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RN 412316-70-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, compd. with hydrazinecarboximidamide (1:1), monoformate (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 79-17-4 CMF C H6 N4

NH || H<sub>2</sub>N-C-NH-NH<sub>2</sub>

CM 3

CRN 64-18-6 CMF C H2 O2

o = CH - OH

RN 412316-71-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-ethoxy-, compd. with hydrazinecarboximidamide (1:1), monoformate (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CM .

CRN 79-17-4 CMF C H6 N4

CM 3

CRN 64-18-6 CMF C H2 O2

o = CH - OH

RN 412316-72-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with hydrazinecarboximidamide (1:1), monoformate (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 79-17-4 CMF C H6 N4

CM 3

CRN 64-18-6 CMF C H2 O2

o = ch - oh

RN 412316-73-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy].alpha.-ethoxy-, compd. with 2-amino-2-(hydroxymethyl)-1,3-propanediol
(1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 77-86-1 CMF C4 H11 N O3

$$\begin{array}{c} {\rm ^{NH_2}} \\ {\rm ^{HO-}\,CH_2-C-CH_2-OH} \\ {\rm ^{CH_2-}\,OH} \end{array}$$

RN 412316-74-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, compd. with 2-amino-2-(hydroxymethyl)-1,3-propanediol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 77-86-1 CMF C4 H11 N O3

$$\begin{array}{c} {\rm NH_2} \\ | \\ {\rm HO-CH_2-C-CH_2-OH} \\ | \\ {\rm CH_2-OH} \end{array}$$

RN 412316-75-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, compd. with 2-amino-2-(hydroxymethyl)-1,3-propanediol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CRN 77-86-1 CMF C4 H11 N O3

$$\begin{array}{c} {\rm NH_2} \\ | \\ {\rm HO-CH_2-C-CH_2-OH} \\ | \\ {\rm CH_2-OH} \end{array}$$

RN 412316-76-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with 2-amino-2-(hydroxymethyl)-1,3-propanediol (1:1) (9CI) (CA INDEX NAME)

CM · 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 77-86-1 CMF C4 H11 N O3

$$\begin{array}{c} {\rm ^{NH_2}} \\ {\rm ^{HO-}\,CH_2-} \\ {\rm ^{C-}\,CH_2^{-}\,OH} \\ {\rm ^{CH_2-}\,OH} \end{array}$$

RN 412316-77-9 CAPLUS

CN D-Glucitol, 1-deoxy-1-(octylamino)-, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 23323-37-7 CMF C14 H31 N O5

Absolute stereochemistry.

RN 412316-78-0 CAPLUS

CN D-Glucitol, 1-deoxy-1-(octylamino)-, (.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0 CMF C21 H25 N O4 S

CRN 23323-37-7 CMF C14 H31 N O5

Absolute stereochemistry.

RN 412316-79-1 CAPLUS

CN D-Glucitol, 1-deoxy-1-(octylamino)-, (.alpha.R)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 23323-37-7 CMF C14 H31 N O5

Absolute stereochemistry.

RN 412316-80-4 CAPLUS

CN D-Glucitol, 1-deoxy-1-(octylamino)-, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CRN 23323-37-7 CMF C14 H31 N O5

Absolute stereochemistry.

RN 412316-81-5 CAPLUS

CN D-Glucitol, 1-deoxy-1-(octylamino)-, (.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 23323-37-7 CMF C14 H31 N O5

Absolute stereochemistry.

$$\begin{array}{c|c}
\text{OH} & \text{OH} & \text{H} \\
\text{R} & \text{R} & \text{S}
\end{array}$$

$$\begin{array}{c|c}
\text{OH} & \text{OH} & \text{OH}
\end{array}$$

RN 412316-82-6 CAPLUS

CN D-Glucitol, 1-deoxy-1-(methylamino)-, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CRN 6284-40-8 CMF C7 H17 N O5

Absolute stereochemistry.

RN 412316-83-7 CAPLUS

CN D-Glucitol, 1-deoxy-1-(methylamino)-, (.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 6284-40-8 CMF C7 H17 N O5

Absolute stereochemistry.

RN 412316-84-8 CAPLUS

CN D-Glucitol, 1-deoxy-1-(methylamino)-, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 6284-40-8 CMF C7 H17 N O5

Absolute stereochemistry.

RN 412316-85-9 CAPLUS

CN D-Glucitol, 1-deoxy-1-(methylamino)-, (.alpha.R)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 6284-40-8 CMF C7 H17 N O5

Absolute stereochemistry.

RN 412317-72-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy].alpha.-ethoxy-, compd. with (.beta.S)-.beta.-aminobenzeneethanol (1:1)
(9CI) (CA INDEX NAME)

CM 1

CRN 223607-24-7 CMF C21 H25 N O4 S

CRN 20989-17-7 CMF C8 H11 N O

Absolute stereochemistry. Rotation (+).

RN 412317-73-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, compd. with (.beta.S)-.beta.-aminobenzeneethanol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-25-8 CMF C21 H25 N O4 S

CRN 20989-17-7 CMF C8 H11 N O

Absolute stereochemistry. Rotation (+).

RN 412317-74-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-ethoxy-, compd. with (.beta.S)-.beta.-aminobenzeneethanol (1:1)
(9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CRN 20989-17-7 CMF C8 H11 N O

Absolute stereochemistry. Rotation (+).

RN 412317-76-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)-, compd. with (.beta.S)-.beta.-aminobenzeneethanol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

CRN 20989-17-7 CMF C8 H11 N O

Absolute stereochemistry. Rotation (+).

RN 412915-55-0 CAPLUS

CN L-Lysine, mono[rel-(+)-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

CRN 56-87-1

CMF C6 H14 N2 O2

Absolute stereochemistry.

RN 412915-57-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)-, compd. with N-cyclohexylcyclohexanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

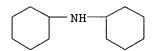
CRN 223607-16-7

CMF C21 H25 N O5

Absolute stereochemistry. Rotation (+).

CM 2

CRN 101-83-7 CMF C12 H23 N



RN 412915-58-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)-, compd. with N,N-dimethylimidodicarbonimidic diamide (1:1) (9CI) (CA INDEX NAME)

CRN 223607-16-7 CMF C21 H25 N O5

Absolute stereochemistry. Rotation (+).

CM 2

CRN 657-24-9 CMF C4 H11 N5

RN 412915-59-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)-, compd. with hydrazinecarboximidamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

CRN 79-17-4 CMF C H6 N4

RN 412915-60-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)-, compd. with (.alpha.R)-.alpha.-methylbenzenemethanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

CRN 3886-69-9 CMF C8 H11 N

Absolute stereochemistry.

RN 412915-63-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)-, formate, compd. with hydrazinecarboximidamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

Absolute stereochemistry. Rotation (+).

CM 2

CRN 79-17-4 CMF C H6 N4

ин || н<sub>2</sub>и-с-ин-ин<sub>2</sub>

CM 3

CRN 64-18-6 CMF C H2 O2 O = CH - OH

RN 412915-66-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)-, compd. with 2-amino-2-(hydroxymethyl)-1,3-propanediol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

Absolute stereochemistry. Rotation (+).

CM 2

CRN 77-86-1 CMF C4 H11 N O3

$$\begin{array}{c} ^{\rm NH_2} \\ | \\ {\rm HO-CH_2-C-CH_2-OH} \\ | \\ {\rm CH_2-OH} \end{array}$$

RN 412915-68-5 CAPLUS

CN D-Glucitol, 1-deoxy-1-(octylamino)-, rel-(+)-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

CRN 23323-37-7 CMF C14 H31 N O5

Absolute stereochemistry.

RN 412915-69-6 CAPLUS

CN D-Glucitol, 1-deoxy-1-(methylamino)-, rel-(+)-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

CRN 6284-40-8 CMF C7 H17 N O5

Absolute stereochemistry.

RN 412915-70-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)-, compd. with (.beta.S)-.beta.-aminobenzeneethanol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 223607-16-7 CMF C21 H25 N O5

CRN 20989-17-7 CMF C8 H11 N O

Absolute stereochemistry. Rotation (+).

RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

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AN
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     133:350232
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     Ranjan, Chakrabarti
     Reddy's Research Foundation, India
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     (substituted) alkyl, cycloalkyl, alkoxy, cycloalkoxy, aryl, aryloxy,
     aralkyl, aralkoxy, heterocyclyl, heteroaryl, amino, acylamino,
     aralkoxycarbonyl, alkylthio, etc.; R5 and/or R6 (on C) may also be oxo; R7
     = H, OH, alkoxy, halo, alkyl, (substituted) aralkyl; R8 = R7, acyl; or
     R7R8 = pi bond; R9 = H, (substituted) alkyl, cycloalkyl, aryl, aralkyl,
     alkoxyalkyl, aryloxycarbonyl, alkylaminocarbonyl, acyl, heterocyclyl,
     heteroaryl, etc.; R10 = H, (substituted) alkyl, cycloalkyl, aryl, aralkyl,
     heterocyclyl, heteroaryl, heteroaralkyl; X = S, O, NR11; R11 = H,
     (substituted) alky, aryl, acyl, etc.; Y = O, NR12; R12 = H, alkyl, aryl,
     hydroxyalkyl, aralkyl, heterocyclyl, heteroaryl, heteroaralkyl; R10R12 may
     form 5- to 6-membered N-contg. heterocyclyl; \bar{n} = 1-4; m = 0, 1; Ar =
     (substituted) single or fused arom. or heterocyclic divalent group]. The
     compds. are agonists at PPAR.alpha. and/or PPAR.gamma., and optionally
     inhibit HMG-CoA reductase. They are thus useful in a variety of
     pharmaceutical roles, particularly as antiobesity and hypocholesterolemic
     agents. Examples include 77 syntheses and 5 precursor prepns., plus
     several bioassays. For instance, title compd. II was prepd. by
     etherification of the corresponding phenol deriv. and benzothiazinylethyl
     mesylate ester (70% yield). At doses of 1-10 mg/kg orally in mice,
     selected compds. I reduced plasma triglyceride levels by 23-74%, blood
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glucose by 45-53%, and total cholesterol levels by 42-57%. IT 223607-15-6P 223607-17-8P, (S)-(-)-2-Ethoxy-3-[4-[2-(2,3dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]propanoic acid 223607-24-7P 223607-27-0P, (S)-(-)-2-Ethoxy-3-[4-[2-(2,3dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]phenyl]propanoic acid 305827-41-2P, (S)-(-)-Ethyl 2-ethoxy-3-[4-[2-(2,3-dihydro-4H-1,4benzothiazin-4-yl)ethoxy]phenyl]propanoate 305827-45-6P, (R)-(+)-Ethyl 2-ethoxy-3-[4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4yl)ethoxy]phenyl]propanoate 305827-61-6P, (S)-(-)-Ethyl 2-ethoxy-3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4yl)ethoxy]phenyl]propanoate RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses) (drug candidate; prepn. of benzoxazines and benzothiazines as antiobesity and hypocholesterolemic agents) RN 223607-15-6 CAPLUS Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-CN .alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-17-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)- (9CI) (CA INDEX NAME)

RN 223607-24-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-27-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)- (9CI) (CA INDEX NAME)

RN 305827-41-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, ethyl ester, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 305827-45-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, ethyl ester, (.alpha.R)- (9CI) (CA INDEX NAME)

RN 305827-61-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, ethyl ester, (.alpha.S)- (9CI) (CA INDEX NAME)

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IT
    223606-56-2P 223606-57-3P 223606-58-4P
    223606-64-2P 223606-65-3P 223606-66-4P
    223606-67-5P 223606-68-6P 223606-70-0P
    223606-71-1P 223606-72-2P 223606-73-3P
    223606-75-5P 223606-76-6P 223606-78-8P
    223606-89-1P 223606-92-6P 223606-95-9P
    223606-97-1P 223606-99-3P 223607-02-1P
    223607-03-2P 223607-04-3P 223607-06-5P
    223607-07-6P 223607-08-7P 223607-16-7P
    223607-25-8P, (R)-(+)-2-Ethoxy-3-[4-[2-(2,3-d)]dihydro-4H-1,4-
    benzothiazin-4-yl)ethoxy]phenyl]propanoic acid 223607-51-0P
    223607-52-1P 223607-54-3P 223607-55-4P
    223607-56-5P 223607-57-6P 223607-58-7P
    223607-59-8P 223607-67-8P 223607-69-0P
    223607-71-4P 223607-73-6P 223607-74-7P
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223607-75-8P 223607-76-9P 223607-77-0P
223607-78-1P 223607-80-5P 223607-82-7P
223607-83-8P 223607-84-9P 223607-86-1P
223607-87-2P 223607-88-3P 223607-89-4P
223607-90-7P 223607-91-8P 223607-92-9P
223607-93-0P 223607-94-1P 223607-95-2P
223607-96-3P 223607-97-4P 223607-99-6P
223608-00-2P 223608-01-3P 223608-02-4P
223608-03-5P 223608-04-6P 223608-05-7P
223608-06-8P 223608-07-9P 223608-08-0P
223608-09-1P 223608-18-2P 223608-19-3P
223608-20-6P 223608-21-7P 223608-22-8P
223608-24-0P 223608-25-1P 223608-26-2P
223608-27-3P 223608-28-4P 305827-50-3P,
3-[4-[2-(2,3-Dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]phenyl]-2-
ethoxypropanoic acid magnesium salt 305827-52-5P,
 (S) - (-) - 3 - [4 - [2 - (2, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (2, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 3 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 3 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 3 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 3 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 3 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4 - yl) ethoxy] et
ethoxypropanoic acid sodium salt 305827-54-7P,
 (S) - (-) - 3 - [4 - [2 - (2, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 -
ethoxypropanoic acid magnesium salt 305827-56-9P,
 (S) - (-) - 3 - [4 - [2 - (2, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (2, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - 2 - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] phenyl] - (3, 3 - Dihydro - 4H - 1, 4 - benzothiazin - 4 - yl) ethoxy] ethoxy
ethoxypropanoic acid arginine salt 305827-58-1P,
3-[4-[2-(2,3-Dihydro-4H-1,4-benzoxazin-4-y1)ethoxy]phenyl]-2-
ethoxypropanoic acid arginine salt 305827-65-0P
305827-67-2P, (S)-(-)-3-[4-[2-(2,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4H-1,4-benzoxazin-4-(3,3-Dihydro-4-(3,3-Dihydro-4-(3,3-Dihydro-4-(3,3-Dihydro-4-(3,3-Dihydro-4-(3,3-Dihydro-4-(3,3-Dih
yl)ethoxy]phenyl]-2-ethoxypropanoic acid arginine salt
305827-89-8P, (+)-Ethyl 3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-
yl)ethoxy]phenyl]-2-hydroxypropanoate 305827-91-2p, (-)-Ethyl
3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-y1)ethoxy]phenyl]-2-
hydroxypropanoate 305827-92-3P, (+)-Ethyl 3-[4-[2-(2,3-dihydro-
4H-1,4-benzothiazin-4-yl)ethoxy]phenyl]-2-hydroxypropanoate
305827-93-4P, (-)-Ethyl 3-[4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-
yl)ethoxy]phenyl]-2-hydroxypropanoate 305827-94-5P, (+)-Ethyl
3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-y1)ethoxy]phenyl]-2-
benzyloxypropanoate 305827-96-7P, (-)-Ethyl 3-[4-[2-(2,3-dihydro-
4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]-2-benzyloxypropanoate
305827-98-9P, (+)-Ethyl 3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-
yl)ethoxy]phenyl]-2-butoxypropanoate 305828-00-6P, (-)-Ethyl
3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-y1)ethoxy]phenyl]-2-
butoxypropanoate 305828-02-8P, (+)-Ethyl 3-[4-[2-(2,3-dihydro-4H-
1,4-benzoxazin-4-yl)ethoxy]phenyl]-2-hexyloxypropanoate
305828-04-0P, (-)-Ethyl 3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-
yl)ethoxy]phenyl]-2-hexyloxypropanoate 305828-20-0P,
(+)-2-Methyl-3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]-2-
ethoxypropanoic acid 305828-22-2p, (-)-2-Methyl-3-[4-[2-(2,3-
dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]-2-ethoxypropanoic acid
305828-24-4P, (+)-2-(2-Fluorobenzyl)-3-[4-[2-(2,3-dihydro-4H-1,4-
benzoxazin-4-yl)ethoxy]phenyl]-2-ethoxypropanoic acid 305828-26-6P
(-) -2-(2-Fluorobenzyl) -3-[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-
yl)ethoxy]phenyl]-2-ethoxypropanoic acid
RL: BAC (Biological activity or effector, except adverse); BSU (Biological
study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
BIOL (Biological study); PREP (Preparation); USES (Uses)
         (drug candidate; prepn. of benzoxazines and benzothiazines as
        antiobesity and hypocholesterolemic agents)
223606-56-2 CAPLUS
Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-
 .alpha.-ethoxy-, methyl ester (9CI) (CA INDEX NAME)
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RN

CN

RN 223606-57-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223606-58-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223606-64-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223606-65-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223606-66-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223606-67-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, methyl ester (9CI) (CA INDEX NAME)

RN 223606-68-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223606-70-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

RN 223606-71-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, methyl ester (9CI) (CA INDEX NAME)

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RN 223606-72-2 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, methyl ester, rel-(+)-(9CI) (CA INDEX NAME)

Rotation (+).

RN 223606-73-3 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, methyl ester, rel-(-)-(9CI) (CA INDEX NAME)

RN 223606-75-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223606-76-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223606-78-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223606-89-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-92-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-95-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-97-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-99-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, ethyl ester (9CI) (CA INDEX NAME)

RN 223607-02-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223607-03-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-04-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-06-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223607-07-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-08-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-16-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 223607-25-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 223607-51-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl- (9CI) (CA INDEX NAME)

RN 223607-52-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro- (9CI) (CA INDEX NAME)

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RN 223607-54-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-55-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-56-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-57-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-58-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-59-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-67-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy- (9CI) (CA INDEX NAME)

RN 223607-69-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-71-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-73-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy- (9CI) (CA INDEX NAME)

RN 223607-74-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-75-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-76-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)- (9CI) (CA INDEX NAME)

RN 223607-77-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-78-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-80-5 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]- (9CI) (CA INDEX NAME)

RN 223607-82-7 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-83-8 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-y1)ethoxy]-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-84-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)- (9CI) (CA INDEX NAME)

RN 223607-86-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-87-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-88-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223607-89-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-90-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-91-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223607-92-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-93-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-94-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223607-95-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-96-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-97-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223607-99-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223608-00-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223608-01-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223608-02-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223608-03-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy].alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX
NAME)

Rotation (-).

RN 223608-04-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223608-05-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223608-06-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223608-07-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, 4-nitrophenyl ester (9CI) (CA INDEX NAME)

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RN 223608-08-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, 4-nitrophenyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223608-09-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, 4-nitrophenyl ester, rel-(-)- (9CI) (CA INDEX NAME)

RN 223608-18-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-19-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-20-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-21-7 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, sodium salt (9CI) (CA INDEX NAME)

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Na

RN 223608-22-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-24-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-25-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-26-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-27-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-28-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 305827-50-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, magnesium salt (9CI) (CA INDEX NAME)

RN 305827-52-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, sodium salt, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

Na

RN 305827-54-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, magnesium salt, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

## ●1/2 Mg

RN 305827-56-9 CAPLUS

CN L-Arginine, mono[(.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CM 1

CRN 223607-27-0 CMF C21 H25 N O4 S

Absolute stereochemistry. Rotation (-).

CM 2

CRN 74-79-3 CMF C6 H14 N4 O2

Absolute stereochemistry.

$$H_2N$$
 $NH$ 
 $NH$ 
 $NH$ 
 $NH$ 
 $NH_2$ 
 $NH_2$ 

RN 305827-58-1 CAPLUS

CN L-Arginine, mono[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CM 1

CRN 223607-15-6 CMF C21 H25 N O5

CM 2

CRN 74-79-3

CMF C6 H14 N4 O2

Absolute stereochemistry.

RN 305827-65-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-

.alpha.-ethoxy-, magnesium salt, (.alpha.S)- (9CI) (CA INDEX NAME)
Absolute stereochemistry. Rotation (-).

## ●1/2 Mg

RN 305827-67-2 CAPLUS

CN L-Arginine, mono[(.alpha.S)-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxybenzenepropanoate] (9CI) (CA INDEX NAME)

CM 1

CRN 223607-17-8 CMF C21 H25 N O5

Absolute stereochemistry. Rotation (-).

CM 2

CRN 74-79-3 CMF C6 H14 N4 O2

Absolute stereochemistry.

RN 305827-89-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester, (+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 305827-91-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 305827-92-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester, (+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 305827-93-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 305827-94-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, ethyl ester, (+)- (9CI) (CA INDEX NAME)

RN 305827-96-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-(phenylmethoxy)-, ethyl ester, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 305827-98-9 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, ethyl ester, (+)- (9CI) (CA INDEX NAME)

RN 305828-00-6 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, ethyl ester, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 305828-02-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, ethyl ester, (+)- (9CI) (CA INDEX NAME)

RN 305828-04-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, ethyl ester, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 305828-20-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, (+)- (9CI) (CA INDEX NAME)

RN 305828-22-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, (-)- (9CI) (CA INDEX NAME)

## Rotation (-).

RN 305828-24-4 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, (+)- (9CI) (CA INDEX NAME)

RN 305828-26-6 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, (-)- (9CI) (CA INDEX NAME)

RE.CNT 12 THERE ARE 12 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

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ANSWER 5 OF 5 CAPLUS COPYRIGHT 2003 ACS on STN
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     Reddy's Research Foundation, India; Reddy-Cheminor, Inc.
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AB
     Title compds. I; R1-R6 = H, halo, OH, NO2, cyano, CHO, (substituted)
     alkyl, cycloalkyl, alkoxy, cycloalkoxy, aryl, aryloxy, aralkyl, aralkoxy,
    heterocyclyl, heteroaryl, amino, acylamino, aralkoxycarbonyl, alkylthio,
     etc.; R7 = H, OH, alkoxy, halo, alkyl, (substituted) aralkyl, bond; R8 =
     R7, acyl; R9 = H, (substituted) alkyl, cycloalkyl, aryl, aralkyl,
     alkoxyalkyl, aryloxycarbonyl, alkylaminocarbonyl, acyl, heterocyclyl,
    heteroaryl, etc.; R10 = H, (substituted) alkyl, cycloalkyl, aryl, aralkyl, heterocyclyl, heteroaryl, heteroaralkyl; Y = O, NR12; R12 = H, alkyl,
     aryl, hydroxyalkyl, aralkyl, heterocyclyl, heteroaryl, heteroaralkyl;
     R10R12 = 5-6 membered (hetero)cyclyl; n = 1-4; m = 0, 1; Ar = 1-4
     (substituted) divalent. Thus, 3-[6-[2-(2,3-dihydro-1,4-benzothiazin-4-
    yl)ethoxy]naphthyl]-2-ethoxypropanoic acid Na salt (prepn. given) at 1
    mg/kg orally in rats reduced plasma triglyceride and total cholesterol
    levels by 50% and 42%, resp.
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223606-64-2P 223606-65-3P 223606-66-4P

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223608-27-3P 223608-28-4P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological
study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
BIOL (Biological study); PREP (Preparation); USES (Uses)
   (prepn. of benzoxazines and benzothiazines as antiobesity and
   hypocholesterolemic agents)
223606-56-2 CAPLUS
Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-
.alpha.-ethoxy-, methyl ester (9CI) (CA INDEX NAME)
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RN

CN

RN 223606-57-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223606-58-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-ethoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223606-64-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223606-65-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223606-66-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

RN 223606-67-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, methyl ester (9CI) (CA INDEX NAME)

RN 223606-68-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223606-70-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-ethoxy-.alpha.-methyl-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223606-71-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, methyl ester (9CI) (CA INDEX NAME)

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RN 223606-72-2 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, methyl ester, rel-(+)-(9CI) (CA INDEX NAME)

RN 223606-73-3 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, methyl ester, rel-(-)-(9CI) (CA INDEX NAME)

Rotation (-).

RN 223606-75-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223606-76-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223606-78-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

RN 223606-89-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-92-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-95-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-97-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, ethyl ester (9CI) (CA INDEX NAME)

RN 223606-99-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, ethyl ester (9CI) (CA INDEX NAME)

RN 223607-02-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223607-03-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-04-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-06-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223607-07-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-08-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-15-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-16-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-ethoxy-, (.alpha.R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 223607-17-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 223607-24-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-25-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 223607-27-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, (.alpha.S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 223607-51-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl- (9CI) (CA INDEX NAME)

RN 223607-52-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-(9CI) (CA INDEX NAME)

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RN 223607-54-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-55-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-56-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-57-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy- (9CI) (CA INDEX NAME)

RN 223607-58-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-59-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, rel-(-)- (9CI) (CA INDEX NAME)

## Rotation (-).

RN 223607-67-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy- (9CI) (CA INDEX NAME)

RN 223607-69-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy].alpha.-hydroxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-71-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-hydroxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-73-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy- (9CI) (CA INDEX NAME)

RN 223607-74-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-75-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-hydroxy-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-76-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)- (9CI) (CA INDEX NAME)

RN 223607-77-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-78-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-80-5 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]- (9CI) (CA INDEX NAME)

RN 223607-82-7 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-83-8 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-84-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)- (9CI) (CA INDEX NAME)

RN 223607-86-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-87-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-88-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223607-89-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-90-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-91-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223607-92-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223607-93-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223607-94-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223607-95-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223607-96-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223607-97-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223607-99-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223608-00-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223608-01-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester (9CI) (CA INDEX NAME)

RN 223608-02-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(+)- (9CI) (CA INDEX NAME)

RN 223608-03-5 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, methyl ester, rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-).

RN 223608-04-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy- (9CI) (CA INDEX NAME)

RN 223608-05-7 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223608-06-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, rel-(-)- (9CI) (CA INDEX NAME)

RN 223608-07-9 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy].alpha.-ethoxy-, 4-nitrophenyl ester (9CI) (CA INDEX NAME)

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RN 223608-08-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, 4-nitrophenyl ester, rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+).

RN 223608-09-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, 4-nitrophenyl ester, rel-(-)- (9CI) (CA INDEX NAME)

RN 223608-18-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-19-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzothiazin-4-yl)ethoxy]-.alpha.-ethoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-20-6 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-.alpha.-methyl-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-21-7 CAPLUS

CN Benzenepropanoic acid, .alpha.-[[4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]phenyl]methyl]-.alpha.-ethoxy-2-fluoro-, sodium salt (9CI) (CA INDEX NAME)

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RN 223608-22-8 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-3-oxo-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-ethoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-24-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(phenylmethoxy)-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-25-1 CAPLUS

CN Benzenepropanoic acid, .alpha.-butoxy-4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-26-2 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-(hexyloxy)-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-27-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-phenoxy-, sodium salt (9CI) (CA INDEX NAME)

RN 223608-28-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-(2,3-dihydro-4H-1,4-benzoxazin-4-yl)ethoxy]-.alpha.-methyl-.alpha.-phenoxy-, sodium salt (9CI) (CA INDEX NAME)

RE.CNT 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

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11 S L1 SSS SAM L2171 S L1 SSS FUL L3

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L45 S L3

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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL

ENTRY SESSION 0.00 -3.26 ENTRY

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